0081_distribution_fit

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1 Distribution fitting to data

SciPy has over 80 distributions that may be used to either generate data or test for fitting of exisitng data. In this example we will test for fit against ten distributions and plot the best three fits. For a full list of distributions see:

https://docs.scipy.org/doc/scipy/reference/stats.html

In this example we'll take the first feature (column) from the Wisconsin Breast Cancer data set. As usual we will start by loading general modules used, and load our data (selecting the first column for our 'y', the data to be fitted).

```
In [12]: import pandas as pd
    import numpy as np
    import scipy
    from sklearn.preprocessing import StandardScaler
    import scipy.stats
    import matplotlib.pyplot as plt
    %matplotlib inline
    # Load data and select first column

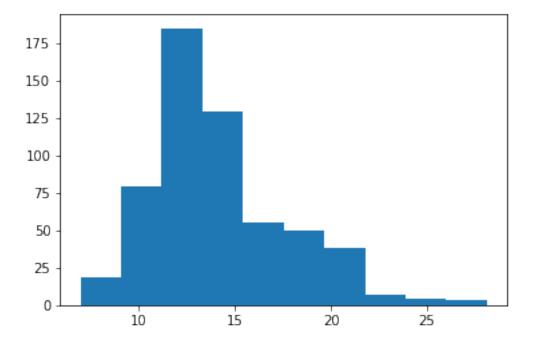
from sklearn import datasets
    data_set = datasets.load_breast_cancer()
    y=data_set.data[:,0]

# Create an index array (x) for data

x = np.arange(len(y))
```

1.1 Visualise the data, and show descriptive summary

Before we do any fitting of distributions, it's always good to do a simple visualisation of the data, and show desriptive statistics. We may, for example, decide to perform some outlier removal if we thinnk that necessary (if there appear to be data points that don't belong to the rest of the population).



If we put the data in a Pandas Dataframe we can use the Pandas secribe method to show a summary.

```
In [14]: y_df = pd.DataFrame(y, columns=['Data'])
         y_df.describe()
Out [14]:
                       Data
                569.000000
         count
         mean
                  14.127292
         std
                   3.524049
                   6.981000
         min
         25%
                  11.700000
         50%
                  13.370000
         75%
                  15.780000
                  28.110000
         max
```

1.2 Fitting a range of distribution and test for goodness of fit

This method will compare will fit a number of distributions to our data, compare goodness of fit with an approximate chi-squared, and test for significant difference between observed and fitted distribution with a Kolmogorov-Smirnov test.

The approximate chi-squared bins data into 50 bins (this could be reduced for smaller data sets) based on percentiles so that each bin contains approximately an equal number of values. For each fitted distribution the expected count of values in each bin predicted from the distribution. The chi-squared value is the the sum of the relative squared error for each bin, such that:

chi-squared = sum ((observed - predicted) ** 2) / predicted)

The lower the chi-squared value the better the fit.

The Kolmogorov-Smirnov test assumes that data has been standardised: that is the mean is subtracted from all data (so the data becomes centred around zero), and that the results values are divided by the standard deviation (so all data becomes expressed as the number of standard deviations above or below the mean). A value of greater than 0.05 means that the fitted distribution is not significantly different to the observed distribution of the data.

It is worth noting that statistical distributions are theoretical models of real-world data. Statistical distributions offer a good way of approximating data (and simplifying huge amounts of data into a few parameters). But when you have a large set of real-world data it is not surprising to find that no theoretical distribution fits the data perfectly. Having he Kolmogorov-Smirnov tests for all distributions produce results of P<0.05 (fitted distribution is statistically different to the observed data distribution) is not unusual for large data sets. In that case in modelling we are generally happy to continue with a fit that looks 'reasonable', being aware this is one of the simplifications present in any model.

Let's first standardise the data using sklearn's StandardScaler:

Now we will fit 10 different distributions, rank them by the approximate chi-squared goodness of fit, and report the Kolmogorov-Smirnov (KS) P value results. Remember that we want chi-squared to be as low as possible, and ideally we want the KS P-value to be >0.05.

Python may report warnings while running the distributions.

```
In [37]: # Set list of distributions to test
         # See https://docs.scipy.org/doc/scipy/reference/stats.html for more
         # Turn off code warnings (this is not recommended for routine use)
         import warnings
         warnings.filterwarnings("ignore")
         # Set up list of candidate distributions to use
         # See https://docs.scipy.org/doc/scipy/reference/stats.html for more
         dist_names = ['beta',
                        'expon',
                        'gamma',
                       'lognorm',
                        'norm',
                       'pearson3',
                        'triang',
                        'uniform',
                        'weibull_min',
                        'weibull_max']
```

```
# Set up empty lists to stroe results
chi_square = []
p_values = []
# Set up 50 bins for chi-square test
# Observed data will be approximately evenly distributed aross all bins
percentile_bins = np.linspace(0,100,51)
percentile_cutoffs = np.percentile(y_std, percentile_bins)
observed_frequency, bins = (np.histogram(y_std, bins=percentile_cutoffs)/
                            sum(y))
# Loop through candidate distributions
for distribution in dist_names:
    # Set up distribution and get fitted distribution parameters
    dist = getattr(scipy.stats, distribution)
    param = dist.fit(y_std)
    # Obtain the KS test P statistic, round it to 5 decimal places
   p = scipy.stats.kstest(y_std, distribution, args=param)[1]
    p = np.around(p, 5)
   p_values.append(p)
    # Get expected counts in percentile bins
    # This is based on a 'cumulative distribution function' (cdf)
    cdf_fitted = dist.cdf(percentile_cutoffs, *param[:-2], loc=param[-2],
                          scale=param[-1])
    expected_frequency = []
    for bin in range(len(percentile_bins)-1):
        expected_cdf_area = cdf_fitted[bin+1] - cdf_fitted[bin]
        expected_frequency.append(expected_cdf_area)
    # calculate chi-squared
    expected frequency = np.array(expected frequency)
    ss = sum (((expected_frequency - observed_frequency) ** 2) / observed_frequency)
    chi_square.append(ss)
# Collate results and sort by goodness of fit (best at top)
results = pd.DataFrame()
results['Distribution'] = dist_names
results['chi_square'] = chi_square
results['p_value'] = p_values
results.sort_values(['chi_square'], inplace=True)
# Report results
```

```
print ('\nDistributions sorted by goodness of fit:')
print ('----')
print (results)
```

Distributions sorted by goodness of fit:

```
Distribution chi_square p_value
      lognorm 13.886012 0.17957
3
0
         beta 14.047573 0.06558
5
     pearson3
              14.065486 0.06152
2
        gamma
              14.065500 0.06151
4
         norm 15.266379 0.00000
6
       triang 18.577091 0.00000
        expon 42.609618 0.00000
1
7
      uniform 57.736913 0.00000
8 weibull_min 71.442730 0.00000
9 weibull_max 307.185921 0.00000
```

We will now take the top three fits, plot the fit and return the sklearn parameters. This time we will fit to the raw data rather than the standardised data.

```
In [39]: # Divide the observed data into 100 bins for plotting (this can be changed)
         number_of_bins = 100
         bin_cutoffs = np.linspace(np.percentile(y,0), np.percentile(y,99),number_of_bins)
         # Create the plot
         h = plt.hist(y, bins = bin_cutoffs, color='0.75')
         # Get the top three distributions from the previous phase
         number_distributions_to_plot = 3
         dist_names = results['Distribution'].iloc[0:number_distributions_to_plot]
         # Create an empty list to stroe fitted distribution parameters
         parameters = []
         # Loop through the distributions of get line fit and paraemters
         for dist_name in dist_names:
             # Set up distribution and store distribution paraemters
             dist = getattr(scipy.stats, dist_name)
             param = dist.fit(y)
             parameters.append(param)
             # Get line for each distribution (and scale to match observed data)
             pdf_fitted = dist.pdf(x, *param[:-2], loc=param[-2], scale=param[-1])
             scale_pdf = np.trapz (h[0], h[1][:-1]) / np.trapz (pdf_fitted, x)
```

```
pdf_fitted *= scale_pdf
    # Add the line to the plot
    plt.plot(pdf_fitted, label=dist_name)
    # Set the plot x axis to contain 99% of the data
    # This can be removed, but sometimes outlier data makes the plot less clear
    plt.xlim(0,np.percentile(y,99))
# Add legend and display plot
plt.legend()
plt.show()
# Store distribution paraemters in a dataframe (this could also be saved)
dist_parameters = pd.DataFrame()
dist_parameters['Distribution'] = (
        results['Distribution'].iloc[0:number_distributions_to_plot])
dist_parameters['Distribution parameters'] = parameters
# Print parameter results
print ('\nDistribution parameters:')
print ('----')
for index, row in dist_parameters.iterrows():
    print ('\nDistribution:', row[0])
    print ('Parameters:', row[1] )
                                                     lognorm
17.5
                                                      beta
                                                      pearson3
15.0
12.5
10.0
 7.5
 5.0
 2.5
 0.0
```

15

20

10

${\tt Distribution\ parameters:}$

Distribution: lognorm

Parameters: (0.3411670333611477, 4.067737189292493, 9.490709944326486)

Distribution: beta

Parameters: (5.250946652630802, 31681.971540166232, 6.175477404020517, 47986.57812196006)

Distribution: pearson3

Parameters: (0.8726704680754525, 14.127306804909308, 3.4698385545042782)